## Amendments to the Claims

Please amend Claim 14. The Claim Listing below will replace all prior versions of the claims in the application:

## **Claim Listing**

1. (Original) A compound represented by the structural formula:

or a pharmaceutically acceptable salt thereof, wherein  $R_1$  and  $R_2$  are independently an aliphatic group, a substituted aliphatic group, an aryl group or a substituted aryl group,

R<sub>10</sub> is -H or unsubstituted alkyl group;

R<sub>6</sub> is a carboxylic acid protecting group; and

Y is a covalent bond or a substituted or unsubstituted straight-chained hydrocarbyl group.

- 2. (Original) The compound of Claim 1 wherein Y is a covalent bond or -C(R<sub>7</sub>R<sub>8</sub>)- and R<sub>7</sub> and R<sub>8</sub> are each independently -H, an aliphatic or substituted aliphatic group, or R<sub>7</sub> is -H and R<sub>8</sub> is a substituted or unsubstituted aryl group, or, R<sub>7</sub> and R<sub>8</sub>, taken together, are a C<sub>2</sub>-C<sub>6</sub> substituted or unsubstituted alkylene group.
- 3. (Original) The compound of Claim 2 wherein  $R_7$  and  $R_8$  are both -H.

- 4. (Original) The compound of Claim 1 wherein R<sub>1</sub> is an aryl group or a substituted aryl group.
- 5. (Original) The compound of Claim 1 wherein R<sub>2</sub> is an alkyl group or a substituted lower alkyl group.
- 6. (Original) The compound of Claim 2 wherein R<sub>2</sub> is methyl or ethyl; R<sub>7</sub> is -H; and R<sub>8</sub> is -H or methyl.
- 7. (Original) The compound of Claim 6 wherein  $R_1$  is phenyl or substituted phenyl.
- 8. (Original) The compound of Claim 7 wherein  $R_1$  is phenyl and  $R_2$  is methyl.
- 9. (Original) The compound of Claim 2 wherein R<sub>1</sub> is an aliphatic group or a substituted aliphatic group.
- 10. (Original) The compound of Claim 2 wherein R<sub>2</sub> is an aliphatic group or a substituted aliphatic group.
- 11. (Original) The compound of Claim 10 wherein R<sub>2</sub> is a lower alkyl group or a substituted lower alkyl group.
- 12. (Original) The compound of Claim 1 wherein  $R_{10}$  is H.
- 13. (Original) The compound of Claim 2 wherein  $R_{10}$  is H.
- 14. (Currently Amended) A compound represented by the structural formula:

$$R_1$$
  $S$   $R_2$   $N$   $R_5$   $R_{10}$ 

or a pharmaceutically acceptable salt thereof, wherein  $R_1$  and  $R_2$  are independently an aliphatic group, a substituted aliphatic group, an aryl group or a substituted aryl group;  $R_5$  is -H or a hydrazine protecting group and  $R_{10}$  is -H or a substituted or unsubstituted alkyl group; wherein  $R_5$  and  $R_{10}$  are not both -H.

- 15. (Original) The compound of Claim 14 wherein  $R_5$  is a hydrazine protecting group when  $R_2$  is an aryl group or a substituted aryl group.
- 16. (Original) The compound of Claim 14 wherein  $R_5$  is -H or a hydrazine protecting group when  $R_2$  is an aliphatic or substituted aliphatic group and  $R_{10}$  is -H or an unsubstituted alkyl group.
- 17. (Original) The compound of Claim 14 wherein R<sub>2</sub> is an aliphatic group or a substituted aliphatic group.
- 18. (Original) The compound of Claim 17 wherein  $R_1$  is an aryl group or a substituted aryl group.
- 19. (Original) The compound of Claim 18 wherein R<sub>2</sub> is an alkyl group or a substituted lower alkyl group.
- 20. (Original) The compound of Claim 19 wherein  $R_2$  is methyl or ethyl.

- 21. (Original) The compound of Claim 14 wherein R<sub>1</sub> is phenyl or substituted phenyl.
- 22. (Original) The compound of Claim 21 wherein  $R_1$  is phenyl and  $R_2$  is methyl.
- 23. (Original) The compound of Claim 21 wherein R<sub>1</sub> is phenyl substituted with one or more groups selected from -OH, -Br, -Cl, -I, -F, -OR<sup>a</sup>, -O-COR<sup>a</sup>, -COR<sup>a</sup>, -CN, -NO<sub>2</sub>, -COOH, -SO<sub>3</sub>H, -NH<sub>2</sub>, -NHR<sup>a</sup>, -N(R<sup>a</sup>R<sup>b</sup>), -COOR<sup>a</sup>, -CHO, -CONH<sub>2</sub>, -CONHR<sup>a</sup>, -NHCOR<sup>a</sup>, -NRCOR<sup>a</sup>, -NHCONH<sub>2</sub>, -NHCONR<sup>a</sup>H, -NHCON(R<sup>a</sup>R<sup>b</sup>), -NR<sup>c</sup>CONH<sub>2</sub>, - $NR^{c}CONR^{a}H$ ,  $-NR^{c}CON(R^{a}R^{b})$ ,  $-C(=NH)-NH_{2}$ ,  $-C(=NH)-NHR^{a}$ ,  $-C(=NH)-N(R^{a}R^{b})$  $C(=NR^c)-NH_2$ ,  $-C(=NR^c)-NHR^a$ ,  $-C(=NR^c)-N(R^aR^b)$ ,  $-NH-C(=NH)-NH_2$ ,  $-NH-C(=NH)-NH_2$  $NHR^a$ , -NH-C(=NH)-N(RaRb), -NH-C(=NRc)-NH<sub>2</sub>, -NH-C(=NR°)-NHR<sup>a</sup>, -NH- $C(=NR^c)-N(R^aR^b)$ ,  $-NrdH-C(=NH)-NH_2$ ,  $-NR^d-C(=NH)-NHR^a$ ,  $-NR^d-C(=NH)-NHR^a$  $N(R^aR^b)$ ,  $-NR^d-C(=NR^c)-NH_2$ ,  $-NR^d-C(=NR^c)-NHR^a$ ,  $-NR^d-C(=NR^c)-N(R^aR^b)$ ,  $-NHNH_2$ ,  $-NR^d-C(=NR^c)-N(R^aR^b)$ ,  $-NR^d-C(=NR^c)-N(R^aR^b)$ ,  $-NHNH_2$ ,  $-NR^d-C(=NR^c)-N(R^aR^b)$ NHNHR<sup>a</sup>, -NHN(R<sup>a</sup>R<sup>b</sup>),-SO<sub>2</sub>NH<sub>2</sub>, -SO<sub>2</sub>NHR<sup>a</sup>, -SO<sub>2</sub>NR<sup>a</sup>R<sup>b</sup>, -CH=CHR<sup>a</sup>, -CH=CR<sup>a</sup>R<sup>b</sup>, -CR<sup>c</sup>=CR<sup>a</sup>R<sup>b</sup>,-CR<sup>c</sup>=CHR<sup>a</sup>, -CR<sup>c</sup>=CR<sup>a</sup>R<sup>b</sup>, -CCR<sup>a</sup>, -SH, -SR<sup>a</sup>, -S(O)<sub>2</sub>R<sup>a</sup>, alkyl groups, substituted alkyl group, non-aromatic heterocyclic group, substituted nonaromatic heterocyclic group, benzyl group, substituted benzyl group, aryl group or substituted aryl group wherein Ra-Rd each independently an alkyl group, substituted alkyl group, benzyl, substituted benzyl, aromatic or substituted aromatic group, or, -N(RaRb), taken together, can also form a substituted or unsubstituted non-aromatic heterocyclic group.
- 24. (Original) The compound of Claim 23, wherein  $R_2$  is methyl.
- 25. (Original) The compound of Claim 14 wherein R<sub>1</sub> is a lower alkyl group and R<sub>2</sub> is a phenyl group substituted with one or more groups selected from -OH, -Br, -Cl, -I, -F, -OR<sup>a</sup>, -O-COR<sup>a</sup>, -COR<sup>a</sup>, -CN, -NO<sub>2</sub>, -COOH, -SO<sub>3</sub>H, -NH<sub>2</sub>, -NHR<sup>a</sup>, -N(R<sup>a</sup>R<sup>b</sup>), -COOR<sup>a</sup>, -CHO, -CONH<sub>2</sub>, -CONHR<sup>a</sup>, -CON(R<sup>a</sup>R<sup>b</sup>), -NHCOR<sup>a</sup>, -NRCOR<sup>a</sup>, -NHCONH<sub>2</sub>, -NHCONR<sup>a</sup>H, -NHCON(R<sup>a</sup>R<sup>b</sup>), -NR<sup>c</sup>CONH<sub>2</sub>, -NR<sup>c</sup>CON(R<sup>a</sup>R<sup>b</sup>),

-C(=NH)-NH<sub>2</sub>, -C(=NH)-NHR<sup>a</sup>, -C(=NH)-N(R<sup>a</sup>R<sup>b</sup>), -C(=NR<sup>c</sup>)-NH<sub>2</sub>, -C(=NR<sup>c</sup>)-NHR<sup>a</sup>, -C(=NR<sup>c</sup>)-N(R<sup>a</sup>R<sup>b</sup>), -NH-C(=NH)-NH<sub>2</sub>, -NH-C(=NH)-NHR<sup>a</sup>, -NH-C(=NH)-N(R<sup>a</sup>R<sup>b</sup>), -NH-C(=NR<sup>c</sup>)-NH<sub>2</sub>, -NH-C(=NR<sup>c</sup>)-NH<sub>2</sub>, -NH-C(=NR<sup>c</sup>)-NH<sub>2</sub>, -NR<sup>d</sup>-C(=NH)-NHR<sup>a</sup>, -NR<sup>d</sup>-C(=NH)-N(R<sup>a</sup>R<sup>b</sup>), -NR<sup>d</sup>-C(=NR<sup>c</sup>)-NH<sub>2</sub>, -NR<sup>d</sup>-C(=NR<sup>c</sup>)-NHR<sup>a</sup>, -NR<sup>d</sup>-C(=NR<sup>c</sup>)-NHR<sup>a</sup>, -NHN(R<sup>a</sup>R<sup>b</sup>), -SO<sub>2</sub>NH<sub>2</sub>, -SO<sub>2</sub>NHR<sup>a</sup>, -SO<sub>2</sub>NR<sup>a</sup>R<sup>b</sup>, -CH=CHR<sup>a</sup>, -CH=CR<sup>a</sup>R<sup>b</sup>, -Crc=CR<sup>a</sup>R<sup>b</sup>, -CR<sup>c</sup>=CHR<sup>a</sup>, -CR<sup>c</sup>=CR<sup>a</sup>R<sup>b</sup>, -CCR<sup>a</sup>, -SH, -SR<sup>a</sup>, -S(O)R<sup>a</sup>, -S(O)<sub>2</sub>R<sup>a</sup>, alkyl groups, substituted alkyl group, non-aromatic heterocyclic group, substituted non-aromatic heterocyclic group, benzyl group, substituted benzyl group, aryl group or substituted aryl group wherein R<sup>a</sup>-R<sup>d</sup> each are independently an alkyl group, substituted alkyl group, benzyl, substituted benzyl, aromatic or substituted aromatic group, or, -N(R<sup>a</sup>R<sup>b</sup>), taken together, can also form a substituted or unsubstituted non-aromatic heterocyclic group.